

**DETAILED ACTION**

Claims 1- 7, 9-21 are pending.

Applicants have elected Group I of the restriction, Claims 1-7, 11-16, 19 and 20 drawn to a formula I where in Y is a N.

***Claim Rejections - 35 USC § 102(e)***

the invention was described in (1) an application for patent, published under section 122(b), by another filed in the United States before the invention by the applicant for patent or (2) a patent granted on an application for patent by another filed in the United States before the invention by the applicant for patent, except that an international application filed under the treaty defined in section 351(a) shall have the effects for purposes of this subsection of an application filed in the United States only if the international application designated the United States and was published under Article 21(2) of such treaty in the English language.

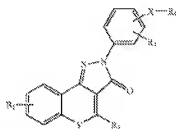
Claims 1-7, 11-16, 19 and 20 are rejected under 35 U.S.C. 102(e) as being anticipated by US 7291612 and US 7081456. Matthews et al .

The applied reference has a common inventor with the instant application. Based upon the earlier effective U.S. filing date of the reference, it constitutes prior art under 35 U.S.C. 102(e). This rejection under 35 U.S.C. 102(e) might be overcome either by a showing under 37 CFR 1.132 that any invention disclosed but not claimed in the reference was derived from the inventor of this application and is thus not the invention "by another," or by an appropriate showing under 37 CFR 1.131.

See claims of the US 7291612

Art Unit: 1625

1. A compound of formula (I) or a pharmaceutically or veterinarily acceptable salt thereof:



(I)

wherein

$R_1$  and  $R_2$  independently represent H; F; Cl; Br;  $-\text{NO}_2$ ;  $-\text{CN}$ ;  $\text{C}_1\text{-C}_6$  alkyl optionally substituted by F or Cl; or  $\text{C}_1\text{-C}_6$  alkoxy optionally substituted by F;  
 $R_3$  represents H, or optionally substituted  $\text{C}_1\text{-C}_6$  alkyl,  $\text{C}_3\text{-C}_7$  cycloalkyl or optionally substituted phenyl;  
 $Y$  represents  $-\text{O}-$ ,  $-\text{S}-$ , N-oxide, or  $-\text{N}(\text{R}_4)-$  wherein  $\text{R}_4$  represents H or  $\text{C}_1\text{-C}_6$  alkyl;  
 $X$  represents a bond or a divalent  $\text{C}_3\text{-C}_6$  alkylene radical;  
 $R_4$  represents  $-\text{NR}_5\text{C}(=\text{O})\text{R}_6$ ,  $-\text{NR}_5\text{C}(=\text{O})\text{OR}_6$ ,  $-\text{NHR}_5(\text{mmS})\text{NHR}_6$  or  $-\text{NHC}(\text{mmS})\text{NHR}_6$  wherein:  
 when  $\text{R}_5$  represents  $-\text{NR}_5\text{C}(=\text{O})\text{R}_6$  or  $-\text{NR}_5\text{C}(=\text{O})\text{OR}_6$ ,  $\text{R}_6$  represents H, or a radical of formula  $-(\text{Alk})_b\text{-Q}$  wherein  $b$  is 0 or 1 and  $\text{Q}$  represents H;  $-\text{CF}_3$ ;  $-\text{OR}$ ;  $-\text{SH}$ ;  $-\text{NR}_8\text{R}_9$  wherein each  $\text{R}_8$  may be the same or different; an ester group; or an optionally substituted phenyl,  $\text{C}_3\text{-C}_7$  cycloalkyl,  $\text{C}_3\text{-C}_7$  cycloalkenyl or heterocyclic ring having from 5 to 8 ring atoms;  
 when  $\text{R}_5$  represents  $-\text{NHC}(\text{mmS})\text{NHR}_6$  or  $-\text{NHC}(\text{mmS})\text{NHR}_6$ ,  $\text{R}_6$  represents a radical of formula  $-(\text{Alk})_b\text{-Q}$  wherein  $b$  is 1 and  $\text{Q}$  represents  $-\text{CF}_3$ ;  $-\text{OR}$ ;  $-\text{SH}$ ;  $-\text{NR}_8\text{R}_9$  wherein each  $\text{R}_8$  may be the same or different; an ester group; or an optionally substituted phenyl,  $\text{C}_3\text{-C}_7$  cycloalkyl,  $\text{C}_3\text{-C}_7$  cycloalkenyl or heterocyclic ring having from 5 to 8 ring atoms;

56

where Alk is a divalent straight chain or branched  $C_1$ - $C_{12}$  alkylate,  $C_1$ - $C_{12}$  alkoxyate or  $C_1$ - $C_{12}$  alkynylate radical which may be interrupted by one or more non-adjacent  $-\text{O}-$ ,  $-\text{S}-$  or  $-\text{N}(\text{R}_9)-$  radicals wherein  $\text{R}_9$  represents H or  $C_1$ - $C_6$  alkyl,  $C_1$ - $C_6$  alkoxy,  $C_1$ - $C_4$  alkynyl, or  $C_1$ - $C_6$  cycloalkyl, and  $\text{R}_9$  represents H or  $C_1$ - $C_6$  alkyl; or when taken together with the atom or atoms to which they are attached  $\text{R}_6$  and  $\text{R}_9$  form an optionally substituted heterocyclic ring having from 5 to 8 ring atoms.

2. A compound as claimed in claim 1 wherein  $\text{R}_1$  is H, F, Cl, methyl or methoxy.

3. A compound as claimed in claim 1 wherein  $\text{R}_2$  is H, methyl, cyclopropyl, phenyl, or fluoro-, chloro-, methyl-, or methoxy-substituted phenyl.

4. A compound as claimed in claim 1 wherein  $\text{R}_3$  is H, F, Cl, methyl, or methoxy.

5. A compound as claimed in claim 1 wherein Y is  $-\text{O}-$ ,  $-\text{S}-$ , or  $-\text{N}(\text{R}_4)-$  wherein  $\text{R}_4$  represents H or methyl.

6. A compound as claimed in claim 1 wherein X is a bond, or a  $-\text{CH}_2-$  or  $-\text{CH}_2\text{CH}_2-$  radical.

7. A compound as claimed in claim 1 wherein  $\text{R}_5$  represents  $-\text{NR}_6\text{C}(=\text{O})\text{R}_6$ ,  $-\text{NR}_6\text{C}(=\text{O})\text{OR}_6$ ,  $-\text{NH}(\text{C}(=\text{O})\text{NR}_6)$  or  $-\text{NH}(\text{C}(=\text{S})\text{NR}_6)$  wherein:

when  $\text{R}_6$  represents  $-\text{NR}_6\text{C}(=\text{O})\text{R}_6$  or  $-\text{NR}_6\text{C}(=\text{O})\text{OR}_6$ ,  $\text{R}_6$  is H or a radical of formula  $-\text{Alk}_p\text{-Q}$  wherein

b is 0 or 1 and

Alk is a  $-(\text{CH}_2)_n-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_3](\text{CH}_2)_p-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_2](\text{CH}_2)_p\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2-$ ,  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$ , or  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$  radical where n is 1, 2, 3 or 4 and m and p are independently 0, 1, 2, 3 or 4, and Q represents H,  $-\text{OH}$ ,  $-\text{COOCH}_3$ , phenyl, cyclopropyl, cyclopentyl, cyclohexyl, pyridyl, furyl, thiocetyl, or oxazolyl, and when  $\text{R}_6$  represents  $-\text{NH}(\text{C}(=\text{O})\text{NR}_6)$  or  $-\text{NH}(\text{C}(=\text{S})\text{NR}_6)$ ,  $\text{R}_6$  is a radical of formula  $-\text{Alk}_p\text{-Q}$  wherein b is 1 and

Alk is a  $-(\text{CH}_2)_n-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_3](\text{CH}_2)_p-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_2](\text{CH}_2)_p\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2-$ ,  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$ , or  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$  radical, where n is 1, 2, 3 or 4 and m and p are independently 0, 1, 2, 3 or 4, and Q represents  $-\text{OH}$ ,  $-\text{COOCH}_3$ , phenyl, cyclopropyl, cyclopentyl, cyclohexyl, pyridyl, furyl, thiocetyl, or oxazolyl, and

$\text{R}_6$  is H, or when taken together with the nitrogen atom to which they are attached  $\text{R}_6$  and  $\text{R}_7$  form a pyrrolidine-2-one or pyrrolidine-2,5-dione ring.

8. A compound as claimed in claim 1 wherein  $\text{R}_1$  is H, F, or Cl,  $\text{R}_2$  is H, F, or Cl,  $\text{R}_3$  is H, F, or Cl, Y is  $-\text{NH}-$ , X is a bond, and  $\text{R}_5$  represents  $-\text{NR}_6\text{C}(=\text{O})\text{R}_6$ ,  $-\text{NR}_6\text{C}(=\text{O})\text{OR}_6$  or  $-\text{NH}(\text{C}(=\text{O})\text{NR}_6)$  wherein:

when  $\text{R}_6$  represents  $-\text{NR}_6\text{C}(=\text{O})\text{R}_6$  or  $-\text{NR}_6\text{C}(=\text{O})\text{OR}_6$ ,  $\text{R}_6$  is H or a radical of formula  $-\text{Alk}_p\text{-Q}$  wherein

b is 0 or 1 and

Alk is a  $-(\text{CH}_2)_n-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_3](\text{CH}_2)_p-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_2](\text{CH}_2)_p\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2-$ ,  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$ , or  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$  radical where n is 1, 2, 3 or 4 and m and p are independently 0, 1, 2, 3 or 4, and Q represents H,  $-\text{OH}$ ,  $-\text{COOCH}_3$ , phenyl, cyclopropyl, cyclopentyl, cyclohexyl, pyridyl, furyl, thiocetyl, or oxazolyl, and when  $\text{R}_6$  represents  $-\text{NH}(\text{C}(=\text{O})\text{NR}_6)$ ,  $\text{R}_6$  is a radical of formula  $-\text{Alk}_p\text{-Q}$  wherein

b is 1 and

Alk is a  $-(\text{CH}_2)_n-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_3](\text{CH}_2)_p-$ ,  $-\text{CH}[(\text{CH}_2)_m\text{CH}_2](\text{CH}_2)_p\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2-$ ,  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$ , or  $-(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-\text{O}-$ ,  $(\text{CH}_2)_n-$  radical, where n is 1, 2, 3 or 4 and m and p are independently 0, 1, 2, 3 or 4, and Q represents H,  $-\text{OH}$ ,  $-\text{COOCH}_3$ , phenyl, cyclopropyl, cyclopentyl, cyclohexyl, pyridyl, furyl, thiocetyl, or oxazolyl, and

Application/Control Number: 10/537,538  
Art Unit: 1625

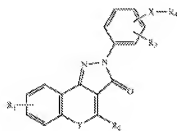
Page 5

And US 7081456.

Art Unit: 1625

FIG. 1

A compound of formula (I) or a pharmaceutically or veterinarily acceptable salt thereof:



(I)

wherein

R<sub>1</sub> and R<sub>2</sub> independently represent H, F, Cl, Br, —NO<sub>2</sub>, —CN, C<sub>1</sub>–C<sub>6</sub> alkyl optionally substituted by F or Cl; or C<sub>1</sub>–C<sub>6</sub> alkoxy optionally substituted by F;

R<sub>3</sub> represents H, or optionally substituted C<sub>1</sub>–C<sub>6</sub> alkyl, C<sub>2</sub>–C<sub>6</sub> cycloalkyl or optionally substituted phenyl;

Y represents —O—, —S—, N-oxide, or —N(R<sub>6</sub>)—, wherein R<sub>6</sub> represents H or C<sub>1</sub>–C<sub>6</sub> alkyl;

X represents a bond or a divalent C<sub>1</sub>–C<sub>6</sub> alkylene radical;

R<sub>4</sub> represents —C(=O)NR<sub>7</sub>R<sub>8</sub>, wherein

R<sub>7</sub> represents a radical of formula —(Alk)<sub>b</sub>—Q wherein b is 1 and

Alk is an optionally substituted divalent straight chain or branched C<sub>1</sub>–C<sub>12</sub> alkylene, C<sub>2</sub>–C<sub>12</sub> alkenylene or C<sub>1</sub>–C<sub>12</sub> alkynylene radical which may be interrupted by one or more non-adjacent —O—, —S— or —N(R<sub>9</sub>)— radicals wherein R<sub>9</sub> represents H or C<sub>1</sub>–C<sub>6</sub> alkyl, C<sub>2</sub>–C<sub>6</sub> alkenyl, C<sub>1</sub>–C<sub>6</sub> alkynyl, or C<sub>2</sub>–C<sub>6</sub> cycloalkyl; and

Q represents H, —CF<sub>3</sub>, —OH, —SH, —NR<sub>10</sub>R<sub>11</sub>, wherein each R<sub>10</sub> may be the same or different; an ester group; or an optionally substituted phenyl, C<sub>2</sub>–C<sub>6</sub> cycloalkyl, C<sub>1</sub>–C<sub>6</sub> cycloalkenyl or heterocyclic ring having from 5 to 8 ring atoms; and

R<sub>5</sub> represents H or C<sub>1</sub>–C<sub>6</sub> alkyl; or when taken together with the atom or atoms to which they are attached R<sub>6</sub> and R<sub>7</sub> form an optionally substituted heterocyclic ring having from 5 to 8 ring atoms.

2. A compound as claimed in claim 1 wherein R<sub>1</sub> is H, F, Cl, methyl or methoxy.

3. A compound as claimed in claim 1 wherein R<sub>2</sub> is H, methyl, methoxy, cyclopropyl, phenyl, or fluoro-, chloro-, methyl, or methoxy-substituted phenyl.

4. A compound as claimed in claim 1 wherein R<sub>3</sub> is H, F, Cl, methyl, or methoxy.

5. A compound as claimed in claim 1 wherein Y is —O—, —S—, or —N(R<sub>6</sub>)— wherein R<sub>6</sub> represents H or methyl.

6. A compound as claimed in claim 1 wherein X is a bond, or a —CH<sub>2</sub>—, or —CH<sub>2</sub>CH<sub>2</sub>— radical.

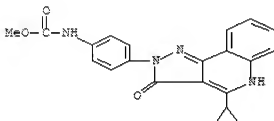
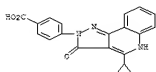
7. A compound as claimed in claim 1 wherein R<sub>4</sub> represents —C(=O)NR<sub>7</sub>R<sub>8</sub>, wherein R<sub>7</sub> is a radical of formula —Alk<sub>b</sub>—Q wherein

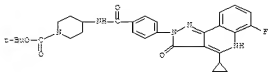
Some of species given are

Art Unit: 1625

RN 702706-05-6 CAPLUS

CN Benzoic acid, 4-(4-cyclopropyl-3,5-dihydro-3-oxo-2H-pyrazolo[4,3-c]quinolin-2-yl)- (CA INDEX NAME)





These compounds read on the applicants when R<sub>2</sub> is a cycloalkyl, R<sub>3</sub> is H, X is a bond.

### ***Double Patenting***

The nonstatutory double patenting rejection is based on a judicially created doctrine grounded in public policy (a policy reflected in the statute) so as to prevent the unjustified or improper timewise extension of the “right to exclude” granted by a patent and to prevent possible harassment by multiple assignees. A nonstatutory obviousness-type double patenting rejection is appropriate where the conflicting claims are not identical, but at least one examined application claim is not patentably distinct from the reference claim(s) because the examined application claim is either anticipated by, or would have been obvious over, the reference claim(s). See, e.g., *In re Berg*, 140 F.3d 1428, 46 USPQ2d 1226 (Fed. Cir. 1998); *In re Goodman*, 11 F.3d 1046, 29 USPQ2d 2010 (Fed. Cir. 1993); *In re Longi*, 759 F.2d 887, 225 USPQ 645 (Fed. Cir. 1985); *In re Van Ornum*, 686 F.2d 937, 214 USPQ 761 (CCPA 1982); *In re Vogel*, 422 F.2d 438, 164 USPQ 619 (CCPA 1970); and *In re Thorington*, 418 F.2d 528, 163 USPQ 644 (CCPA 1969).

A timely filed terminal disclaimer in compliance with 37 CFR 1.321(c) or 1.321(d) may be used to overcome an actual or provisional rejection based on a nonstatutory double patenting ground provided the conflicting application or patent either is shown to be commonly owned with this application, or claims an invention made as a result of activities undertaken within the scope of a joint research agreement.

Effective January 1, 1994, a registered attorney or agent of record may sign a terminal disclaimer. A terminal disclaimer signed by the assignee must fully comply with 37 CFR 3.73(b).

Claims 1-7, 11-16, 19 and 20 are rejected on the ground of nonstatutory double patenting over claims 1-8 of U. S. Patent No. 7291612, and claims 1-7 of US 7081456.

Claims 1-7, 11-16, 19 and 20 are rejected on the ground of nonstatutory obviousness-type double patenting as being unpatentable over claims 1-8, of U.S. Patent No. US 7291612 and claims 1-7 of US Patent No. 7081456.. Although the conflicting claims are not identical, they are not patentably distinct from each other because there is overlapping subject matter.

See rejection above.

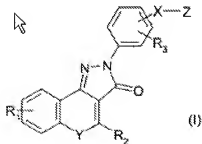
***Claim Rejections - 35 USC § 103***

The following is a quotation of 35 U.S.C. 103(a) which forms the basis for all obviousness rejections set forth in this Office action:

(a) A patent may not be obtained though the invention is not identically disclosed or described as set forth in section 102 of this title, if the differences between the subject matter sought to be patented and the prior art are such that the subject matter as a whole would have been obvious at the time the invention was made to a person having ordinary skill in the art to which said subject matter pertains. Patentability shall not be negated by the manner in which the invention was made.

Claims 1-7, 11-16, 19 and 20 are rejected under 35 U.S.C. 103(a) as being unpatentable over WO 03/004495 (filing date) Bjork et al. US 6642249. (July 2001)

Applicants claims are drawn to compounds and pharmaceutical compositions of the formula



(I)

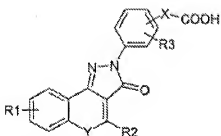
wherein Z is a COOH or an ester

thereof., R<sub>2</sub> is an optionally substituted cycloalkyl or a phenyl



*Scope & Content of Prior Art MPEP 2141.01*

WO 03/004495 Bjork et al teaches compounds of the formula



(I)

Y is a NR<sub>4</sub>, R<sub>2</sub> is a H or a lower alkyl. (

wherein lower alkyl meaning includes cyclic alkyl groups having 1-6 carbon atoms ( see line 9 and 10 page 4 of the reference.

*Difference between Prior Art and the claims MPEP 2141.02*

Even though WO '495 Bjork et al teaches the same core of the same use, it generically teaches R<sub>2</sub> being a cycloalkyl, no species have been made.

*Prima Facie Obviousness , Rational and Motivation MPEP 2142-2413*

Bjork et al generically teaches the same core with Z being a COOH or its ester.

Even though no species have been made the teaching that a cycloalkyl ring can be present at the R<sub>2</sub> substitution and still retain the properties. Thus motivating a person of skill in the art to substitute a carbocyclic ring for R<sub>2</sub> ( cycloalkyl or a phenyl ) and still have some

expectation of success that the compounds would have activity. KSR International v Telflex Inc.

***Conclusion***

Claims 1-7, 11-16, 19 and 20 are not allowable.

Any inquiry concerning this communication or earlier communications from the examiner should be directed to Rita J. Desai whose telephone number is 571-272-0684. The examiner can normally be reached on Monday - Friday, flex time..

If attempts to reach the examiner by telephone are unsuccessful, the examiner's supervisor, Janet Andres can be reached on 571-272-0867. The fax phone number for the organization where this application or proceeding is assigned is 571-273-8300.

Information regarding the status of an application may be obtained from the Patent Application Information Retrieval (PAIR) system. Status information for published applications may be obtained from either Private PAIR or Public PAIR. Status information for unpublished applications is available through Private PAIR only. For more information about the PAIR system, see <http://pair-direct.uspto.gov>. Should you have questions on access to the Private PAIR system, contact the Electronic Business Center (EBC) at 866-217-9197 (toll-free). If you would like assistance from a USPTO Customer Service Representative or access to the automated information system, call 800-786-9199 (IN USA OR CANADA) or 571-272-1000.

/Rita J. Desai/  
Primary Examiner, Art Unit 1625

R.D.  
September 22, 2008

